

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP111824\
 Data File : BP023209.D
 Acq On : 18 Nov 2024 22:47
 Operator : RC/JU
 Sample : SP6687
 Misc : SFAM SURROGATE
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SP6687

Manual Integrations
 APPROVED

Reviewed By :Yogesh Patel 11/19/2024
 Supervised By :mohammad ahmed 11/19/2024

Quant Time: Nov 19 02:27:30 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP110924.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Nov 13 04:49:01 2024
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.581	152	71081	20.000	ng/u1	0.00
20) Naphthalene-d8	10.316	136	272019	20.000	ng/u1	-0.01
38) Acenaphthene-d10	14.181	164	201331	20.000	ng/u1	-0.02
64) Phenanthrene-d10	16.998	188	499922	20.000	ng/u1	#-0.02
79) Chrysene-d12	21.433	240	494789	20.000	ng/u1	-0.01
88) Perylene-d12	24.651	264	543995	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.187	96	29798m	20.150	ng/uL	0.05
4) Pyridine-d5	3.593	84	432681m	98.482	ng/u1	0.05
7) Phenol-d5	6.787	99	538748	102.517	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	6.928	67	298067	96.538	ng/u1	0.00
11) 2-Chlorophenol-d4	7.128	132	400586	104.081	ng/u1	0.00
15) 4-Methylphenol-d8	8.287	113	428487	99.082	ng/u1	0.00
21) Nitrobenzene-d5	8.722	128	182353	96.293	ng/u1	0.00
24) 2-Nitrophenol-d4	9.428	143	219835	97.373	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	9.963	165	467775	98.938	ng/u1	0.00
31) 4-Chloroaniline-d4	10.469	131	471914	83.666	ng/u1	-0.01
46) Dimethylphthalate-d6	13.604	166	1437483	99.065	ng/u1	0.00
49) Acenaphthylene-d8	13.875	160	1490136	98.419	ng/u1	-0.02
54) 4-Nitrophenol-d4	14.457	143	180118m	82.598	ng/u1	0.00
60) Fluorene-d10	15.204	176	1224398	96.242	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.351	200	232310	77.572	ng/u1	-0.01
73) Anthracene-d10	17.092	188	2056037	97.133	ng/u1	-0.02
81) Pyrene-d10	19.492	212	2453231	107.643	ng/u1	0.00
92) Benzo(a)pyrene-d12	24.427	264	2585641	99.783	ng/u1	-0.02

Target Compounds Qvalue

(#) = qualifier out of range (m) = manual integration (+) = signals summed

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