

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP121922\
 Data File : BP013232.D
 Acq On : 22 Dec 2022 05:25
 Operator : CG/JU
 Sample : PB149716BL
 Misc :
 ALS Vial : 115 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK716

Quant Time: Dec 22 06:34:33 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP120722.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Dec 21 23:38:37 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.946	152	420974	20.000	ng/u1	0.00
20) Naphthalene-d8	10.757	136	1634237	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.587	164	942683	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.345	188	1985362	20.000	ng/u1	0.00
79) Chrysene-d12	21.433	240	2096306	20.000	ng/u1	0.00
88) Perylene-d12	23.910	264	2403925	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.334	96	71514	7.263	ng/uL	0.00
4) Pyridine-d5	3.758	84	1005839	35.963	ng/u1	0.00
7) Phenol-d5	7.105	99	1193620	34.810	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.275	67	767115	37.086	ng/u1	0.00
11) 2-Chlorophenol-d4	7.475	132	981577	36.364	ng/u1	0.00
15) 4-Methylphenol-d8	8.658	113	921022	32.782	ng/u1	0.00
21) Nitrobenzene-d5	9.116	128	453445	37.765	ng/u1	0.00
24) 2-Nitrophenol-d4	9.840	143	506748	37.488	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.381	165	928691	35.370	ng/u1	0.00
31) 4-Chloroaniline-d4	10.899	131	1251737	33.770	ng/u1	0.00
46) Dimethylphthalate-d6	13.998	166	2564148	35.392	ng/u1	0.00
49) Acenaphthylene-d8	14.281	160	2942092	36.956	ng/u1	0.00
54) 4-Nitrophenol-d4	14.787	143	369141	28.325	ng/u1	0.00
60) Fluorene-d10	15.581	176	2230206	36.386	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.704	200	320479	25.084	ng/u1	0.00
73) Anthracene-d10	17.445	188	3593223	40.117	ng/u1	0.00
81) Pyrene-d10	19.675	212	4582300	38.082	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.751	264	4865802	40.609	ng/u1	0.00

Target Compounds Qvalue

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

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