

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP111423\
 Data File : BP018152.D
 Acq On : 14 Nov 2023 12:47
 Operator : MA/JU
 Sample : PB156922BL
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK922

Quant Time: Nov 15 00:37:40 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP111023.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Mon Nov 13 15:50:18 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.787	152	66034	20.000	ng/u1	0.00
20) Naphthalene-d8	10.605	136	269643	20.000	ng/u1	# 0.00
38) Acenaphthene-d10	14.469	164	149584	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.251	188	365592	20.000	ng/u1	0.00
79) Chrysene-d12	21.380	240	327984	20.000	ng/u1	#-0.01
88) Perylene-d12	23.851	264	361463	20.000	ng/u1	-0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.158	96	13208	7.056	ng/uL	0.00
4) Pyridine-d5	3.593	84	115087	23.580	ng/u1	-0.01
7) Phenol-d5	6.958	99	196363	30.801	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.140	67	130715	34.890	ng/u1	0.00
11) 2-Chlorophenol-d4	7.311	132	157636	31.225	ng/u1	0.00
15) 4-Methylphenol-d8	8.510	113	154384	29.553	ng/u1	0.00
21) Nitrobenzene-d5	8.999	128	77055	31.719	ng/u1	0.00
24) 2-Nitrophenol-d4	9.710	143	82339	29.917	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.234	165	132316	26.974	ng/u1	0.00
31) 4-Chloroaniline-d4	10.793	131	182716	27.162	ng/u1	0.00
46) Dimethylphthalate-d6	13.893	166	502916	36.438	ng/u1	0.00
49) Acenaphthylene-d8	14.169	160	507062	33.731	ng/u1	0.00
54) 4-Nitrophenol-d4	14.745	143	37305	18.088	ng/u1	0.00
60) Fluorene-d10	15.469	176	421945	37.029	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.634	200	44135	16.699	ng/u1	-0.03
73) Anthracene-d10	17.351	188	679565	34.369	ng/u1	0.00
81) Pyrene-d10	19.604	212	791417	36.125	ng/u1	-0.01
92) Benzo(a)pyrene-d12	23.686	264	735577	34.997	ng/u1	-0.02

Target Compounds Qvalue

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

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