

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP113022\
 Data File : BP012876.D
 Acq On : 30 Nov 2022 13:14
 Operator : CG/JU
 Sample : PB149338BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK338

Quant Time: Nov 30 21:54:35 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP112522.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Nov 29 22:53:34 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.010	152	515915	20.000	ng/u1	0.00
20) Naphthalene-d8	10.828	136	2394759	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.645	164	1690947	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.398	188	3758100	20.000	ng/u1	0.00
79) Chrysene-d12	21.486	240	3397882	20.000	ng/u1	0.00
88) Perylene-d12	23.998	264	2995504	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.393	96	93456	7.762	ng/uL	0.00
4) Pyridine-d5	3.816	84	1228053	34.084	ng/u1	0.00
7) Phenol-d5	7.157	99	1659601	39.147	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.334	67	1045844	40.035	ng/u1	0.00
11) 2-Chlorophenol-d4	7.534	132	1327706	41.101	ng/u1	0.00
15) 4-Methylphenol-d8	8.710	113	1385063	40.550	ng/u1	0.00
21) Nitrobenzene-d5	9.175	128	670032	45.713	ng/u1	0.00
24) 2-Nitrophenol-d4	9.904	143	767138	49.292	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.440	165	1429135	39.355	ng/u1	0.00
31) 4-Chloroaniline-d4	10.957	131	1999922	40.218	ng/u1	0.00
46) Dimethylphthalate-d6	14.051	166	5101730	43.833	ng/u1	0.00
49) Acenaphthylene-d8	14.339	160	5506969	40.551	ng/u1	0.00
54) 4-Nitrophenol-d4	14.828	143	807729	38.925	ng/u1	0.00
60) Fluorene-d10	15.639	176	4406020	43.225	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.751	200	735455	40.945	ng/u1	0.00
73) Anthracene-d10	17.498	188	7376958	44.593	ng/u1	0.00
81) Pyrene-d10	19.727	212	8425253	42.673	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.839	264	6859480	46.620	ng/u1	0.00

Target Compounds Qvalue

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

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