

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP111723\
 Data File : BP018229.D
 Acq On : 17 Nov 2023 20:15
 Operator : MA/JU
 Sample : PB157215BL
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :

Quant Time: Nov 17 21:53:48 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP111023.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Nov 15 18:13:37 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.781	152	42081	20.000	ng/ul	0.00
20) Naphthalene-d8	10.599	136	148309	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.463	164	90020	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.245	188	201260	20.000	ng/ul	0.00
79) Chrysene-d12	21.380	240	179550	20.000	ng/ul	0.00
88) Perylene-d12	23.851	264	191493	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.158	96	11602	9.726	ng/uL	0.00
4) Pyridine-d5	3.593	84	129194	41.537	ng/ul	0.00
7) Phenol-d5	6.952	99	152677	37.581	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.128	67	97716	40.928	ng/ul	0.00
11) 2-Chlorophenol-d4	7.299	132	133212	41.406	ng/ul	0.00
15) 4-Methylphenol-d8	8.505	113	130147	39.094	ng/ul	0.00
21) Nitrobenzene-d5	8.987	128	63118	47.238	ng/ul	0.00
24) 2-Nitrophenol-d4	9.705	143	71844	47.460	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.228	165	114231	42.339	ng/ul	0.00
31) 4-Chloroaniline-d4	10.781	131	156607	42.327	ng/ul	0.00
46) Dimethylphthalate-d6	13.887	166	441690	53.178	ng/ul	0.00
49) Acenaphthylene-d8	14.157	160	451406	49.898	ng/ul	0.00
54) 4-Nitrophenol-d4	14.745	143	35817	28.858	ng/ul	0.01
60) Fluorene-d10	15.463	176	358431	52.269	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.716	200	845	0.581	ng/ul	0.09
73) Anthracene-d10	17.345	188	570661	52.427	ng/ul	0.00
81) Pyrene-d10	19.598	212	684646	57.086	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.680	264	626344	56.251	ng/ul	0.00

Target Compounds Qvalue

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

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