

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP090723\
 Data File : BP016978.D
 Acq On : 07 Sep 2023 09:28
 Operator : MA/JU
 Sample : PB155165BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK165

Quant Time: Sep 08 00:37:03 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP090523.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Sep 06 01:52:05 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.981	152	369557	20.000	ng/u1	0.00
20) Naphthalene-d8	10.804	136	1593719	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.634	164	994389	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.404	188	2135987	20.000	ng/u1	0.00
79) Chrysene-d12	21.504	240	1660973	20.000	ng/u1	0.00
88) Perylene-d12	24.045	264	1589871	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.328	96	79566	8.144	ng/uL	0.00
4) Pyridine-d5	3.764	84	971771	34.810	ng/u1	0.00
7) Phenol-d5	7.128	99	1142971	33.853	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.316	67	748147	35.426	ng/u1	0.00
11) 2-Chlorophenol-d4	7.499	132	861423	35.253	ng/u1	0.00
15) 4-Methylphenol-d8	8.687	113	892049	34.681	ng/u1	0.00
21) Nitrobenzene-d5	9.175	128	436783	35.326	ng/u1	0.00
24) 2-Nitrophenol-d4	9.893	143	481452	35.921	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.416	165	820564	33.162	ng/u1	0.00
31) 4-Chloroaniline-d4	10.969	131	1244197	34.851	ng/u1	0.00
46) Dimethylphthalate-d6	14.051	166	2743786	37.068	ng/u1	0.00
49) Acenaphthylene-d8	14.334	160	3052028	35.728	ng/u1	0.00
54) 4-Nitrophenol-d4	14.851	143	478579	33.100	ng/u1	0.00
60) Fluorene-d10	15.634	176	2314362	36.519	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.757	200	334666	26.035	ng/u1	0.00
73) Anthracene-d10	17.504	188	3652473	37.235	ng/u1	0.00
81) Pyrene-d10	19.745	212	3966349	39.205	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.880	264	3142439	38.960	ng/u1	0.00

Target Compounds Qvalue

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

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