

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP121022\
 Data File : BP013041.D
 Acq On : 10 Dec 2022 15:53
 Operator : CG/JU
 Sample : PB149500BL
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK500

Quant Time: Dec 12 03:25:12 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP120722.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Mon Dec 12 01:32:54 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.969	152	463357	20.000	ng/u1	0.00
20) Naphthalene-d8	10.787	136	2026055	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.610	164	1354074	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.369	188	3032957	20.000	ng/u1	0.00
79) Chrysene-d12	21.451	240	2434805	20.000	ng/u1	0.00
88) Perylene-d12	23.939	264	2234089	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.352	96	85753	7.913	ng/uL	0.00
4) Pyridine-d5	3.781	84	1200810	39.007	ng/u1	0.00
7) Phenol-d5	7.128	99	1581995	41.917	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.299	67	984191	43.229	ng/u1	0.00
11) 2-Chlorophenol-d4	7.499	132	1264511	42.561	ng/u1	0.00
15) 4-Methylphenol-d8	8.675	113	1285724	41.576	ng/u1	0.00
21) Nitrobenzene-d5	9.140	128	648329	43.553	ng/u1	0.00
24) 2-Nitrophenol-d4	9.863	143	751039	44.815	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.399	165	1340698	41.187	ng/u1	0.00
31) 4-Chloroaniline-d4	10.922	131	1834258	39.915	ng/u1	0.00
46) Dimethylphthalate-d6	14.016	166	4378937	42.078	ng/u1	0.00
49) Acenaphthylene-d8	14.304	160	4907249	42.914	ng/u1	0.00
54) 4-Nitrophenol-d4	14.804	143	714308	38.158	ng/u1	0.00
60) Fluorene-d10	15.604	176	3813069	43.310	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.722	200	656005	33.611	ng/u1	0.00
73) Anthracene-d10	17.469	188	6169089	45.086	ng/u1	0.00
81) Pyrene-d10	19.698	212	6802648	48.675	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.780	264	5091597	45.724	ng/u1	0.00

Target Compounds Qvalue

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

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