

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP121022\
 Data File : BP013059.D
 Acq On : 11 Dec 2022 02:41
 Operator : CG/JU
 Sample : PB149488BL
 Misc :
 ALS Vial : 30 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :

Quant Time: Dec 12 03:30:59 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	662996	20.000	ng/ul	0.00
20) Naphthalene-d8	10.781	136	2913932	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.610	164	2000372	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.363	188	4583917	20.000	ng/ul	0.00
79) Chrysene-d12	21.451	240	4203368	20.000	ng/ul	0.00
88) Perylene-d12	23.939	264	4160374	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.352	96	93375	6.022	ng/uL	0.00
4) Pyridine-d5	3.775	84	1269763	28.827	ng/ul	0.00
7) Phenol-d5	7.122	99	1686308	31.226	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.293	67	1070786	32.870	ng/ul	0.00
11) 2-Chlorophenol-d4	7.498	132	1390407	32.707	ng/ul	0.00
15) 4-Methylphenol-d8	8.675	113	1379212	31.170	ng/ul	0.00
21) Nitrobenzene-d5	9.140	128	707722	33.057	ng/ul	0.00
24) 2-Nitrophenol-d4	9.863	143	817403	33.913	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.398	165	1441038	30.781	ng/ul	0.00
31) 4-Chloroaniline-d4	10.922	131	1949652	29.499	ng/ul	0.00
46) Dimethylphthalate-d6	14.016	166	5008723	32.580	ng/ul	0.00
49) Acenaphthylene-d8	14.304	160	5423608	32.105	ng/ul	0.00
54) 4-Nitrophenol-d4	14.798	143	798237	28.865	ng/ul	0.00
60) Fluorene-d10	15.604	176	4237427	32.580	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.722	200	813748	27.586	ng/ul	0.00
73) Anthracene-d10	17.463	188	6867673	33.209	ng/ul	0.00
81) Pyrene-d10	19.692	212	8115752	33.637	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.774	264	6979707	33.658	ng/ul	0.00

Target Compounds	Qvalue

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

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