

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM110222\
 Data File : BM037320.D
 Acq On : 02 Nov 2022 17:29
 Operator : CG/JU
 Sample : PB148669BL
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SBLK669

Quant Time: Nov 02 22:54:51 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM110122.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Nov 02 03:33:37 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.857	152	352074	20.000	ng/u1	0.00
20) Naphthalene-d8	10.657	136	1539925	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.492	164	1007350	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.233	188	2072260	20.000	ng/u1	0.00
79) Chrysene-d12	21.403	240	1856022	20.000	ng/u1	0.00
88) Perylene-d12	23.750	264	1615247	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.275	96	45614	5.711	ng/uL	0.00
4) Pyridine-d5	3.698	84	637648	26.745	ng/u1	0.00
7) Phenol-d5	7.028	99	790767	25.808	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.192	67	520907	27.705	ng/u1	0.00
11) 2-Chlorophenol-d4	7.386	132	657551	28.088	ng/u1	0.00
15) 4-Methylphenol-d8	8.575	113	646423	25.836	ng/u1	0.00
21) Nitrobenzene-d5	9.027	128	321620	27.960	ng/u1	0.00
24) 2-Nitrophenol-d4	9.745	143	350959	28.009	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.280	165	611683	26.478	ng/u1	0.00
31) 4-Chloroaniline-d4	10.804	131	939627	26.617	ng/u1	0.00
46) Dimethylphthalate-d6	13.910	166	2038941	29.274	ng/u1	0.00
49) Acenaphthylene-d8	14.186	160	2363115	28.221	ng/u1	0.00
54) 4-Nitrophenol-d4	14.704	143	315622	22.671	ng/u1	0.00
60) Fluorene-d10	15.480	176	1887433	30.388	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.609	200	295946	22.548	ng/u1	0.00
73) Anthracene-d10	17.327	188	2937568	30.780	ng/u1	0.00
81) Pyrene-d10	19.621	212	3212739	33.125	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.603	264	2569896	30.908	ng/u1	0.00

Target Compounds Qvalue

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

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