

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP102121\
 Data File : BP007599.D
 Acq On : 22 Oct 2021 20:55
 Operator : CG/JU
 Sample : PB140091BL
 Misc :
 ALS Vial : 59 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK091

Quant Time: Oct 23 00:21:40 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SFAM-EPA-BP101421.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Oct 20 10:49:33 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.946	152	194079	20.000	ng/u1	0.00
20) Naphthalene-d8	10.752	136	788783	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.581	164	518482	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.334	188	1125245	20.000	ng/u1	0.00
79) Chrysene-d12	21.422	240	1196838	20.000	ng/u1	0.00
88) Perylene-d12	23.863	264	1224474	20.000	ng/u1	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.376	96	32318	6.577	ng/uL	0.00
4) Pyridine-d5	3.787	84	478301	35.558	ng/u1	0.00
7) Phenol-d5	7.111	99	625585	41.275	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.275	67	334139	36.963	ng/u1	0.00
11) 2-Chlorophenol-d4	7.475	132	492003	41.264	ng/u1	0.00
15) 4-Methylphenol-d8	8.652	113	495821	42.470	ng/u1	0.00
21) Nitrobenzene-d5	9.105	128	232793	46.635	ng/u1	0.00
24) 2-Nitrophenol-d4	9.828	143	259999	55.762	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.369	165	495577	43.576	ng/u1	0.00
31) 4-Chloroaniline-d4	10.881	131	655229	42.605	ng/u1	0.00
46) Dimethylphthalate-d6	13.993	166	1499618	41.806	ng/u1	0.00
49) Acenaphthylene-d8	14.275	160	1783311	39.559	ng/u1	0.00
54) 4-Nitrophenol-d4	14.769	143	259501	45.086	ng/u1	0.00
60) Fluorene-d10	15.575	176	1266712	42.198	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.687	200	56740	12.584	ng/u1	0.00
73) Anthracene-d10	17.434	188	2098474	43.913	ng/u1	0.00
81) Pyrene-d10	19.669	212	2488115	41.585	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.710	264	2562679	42.061	ng/u1	0.00

Target Compounds Qvalue

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

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